

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089965.D
 Acq On : 27 Jun 2015 8:09
 Operator : TP/IZ
 Sample : G2782-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X3SS0230006-150624

Manual Integrations
 APPROVED

apatel
 6/29/2015 5:03:24 PM

Quant Time: Jun 29 01:15:55 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 16:51:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	23042	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	109262	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	66743	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	156017	5.00	ng	0.00
25) Chrysene-d12	21.13	240	193109	5.00	ng	0.00
32) Perylene-d12	23.47	264	176833	5.00	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.27	112	45564	10.95	ng	0.00
5) Phenol-d6	6.81	99	74832	11.74	ng	0.00
7) Nitrobenzene-d5	8.77	82	55126	6.78	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	10860	9.18	ng	0.00
14) 2-Fluorobiphenyl	12.87	172	126855	6.52	ng	0.00
27) Terphenyl-d14	19.60	244	202243	6.07	ng	0.00

Target Compounds

						Qvalue
8) Nitrobenzene	8.81	77	868	0.10	ng	# 97
9) Naphthalene	10.44	128	2850	0.12	ng	# 92
10) Hexachlorobutadiene	10.74	225	258	0.07	ng	97
11) 2-Methylnaphthalene	12.05	142	2196	0.13	ng	96
15) Acenaphthylene	13.95	152	17364	0.15	ng	99
16) Acenaphthene	14.30	154	2986	0.17	ng	91
17) Fluorene	15.29	166	3817	0.17	ng	# 95
19) 4-Bromophenyl-phenylether	16.18	248	933	0.15	ng	94
20) Hexachlorobenzene	16.30	284	4210	0.15	ng	93
21) Pentachlorophenol	16.63	266	1819	1.56	ng	90
22) Phenanthrene	17.01	178	16672	0.43	ng	99
23) Anthracene	17.12	178	7872	0.22	ng	98
24) Fluoranthene	19.02	202	37963	0.89	ng	100
26) Pyrene	19.38	202	37493	0.77	ng	97
28) Benzo(a)anthracene	21.12	228	31328	0.68	ng	98
29) Chrysene	21.17	228	34003	0.67	ng	96
30) Bis(2-ethylhexyl)phthalate	21.07	149	5542	1.41	ng	# 85
31) Indeno(1,2,3-cd)pyrene	25.88	276	20393	0.55	ng	# 100
33) Benzo(b)fluoranthene	22.76	252	37136	0.99	ng	96
34) Benzo(k)fluoranthene	22.80	252	21520m	0.51	ng	
35) Benzo(a)pyrene	23.36	252	26176	0.79	ng	98
36) Dibenzo(a,h)anthracene	25.90	278	9058	0.29	ng	# 91
37) Benzo(g,h,i)perylene	26.62	276	19682	0.59	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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